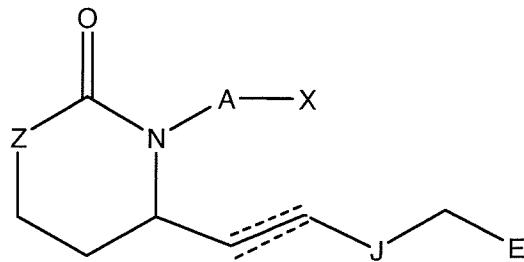


**Amendments to the Claims:**

This listing of claims will replace all prior versions, and listings, of claims in the application:

**Listing of Claims:**

CLAIM 1. (Currently Amended) A compound comprising represented by the formula:



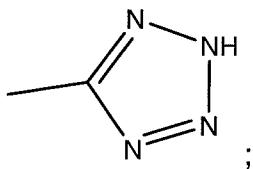
or a pharmaceutically acceptable salt or a prodrug thereof;

wherein a dashed line represents the presence or absence of a double bond or a triple bond;

A is  $-(\text{CH}_2)_6-$ , *cis*  $-\text{CH}_2\text{CH}=\text{CH}-(\text{CH}_2)_3-$ , or  $-\text{CH}_2\text{C}\equiv\text{C}-(\text{CH}_2)_3-$ , wherein 1 or 2 carbon atoms may be substituted with S or O;

Z is O, S, or NR;

X is selected from the group consisting of  $\text{CO}_2\text{H}$ ,  $\text{CONHR}_2$ ,  $\text{CONR}_2$ ,  $\text{CON}(\text{OR})\text{R}$ ,  $\text{CON}(\text{CH}_2\text{CH}_2\text{OH})_2$ ,  $\text{CONH}(\text{CH}_2\text{CH}_2\text{OH})$ ,  $\text{CH}_2\text{OH}$ ,  $\text{P}(\text{O})(\text{OH})_2$ ,  $\text{CONHSO}_2\text{R}$ ,  $\text{SO}_2\text{NR}_2$ ,  $\text{SO}_2\text{NHR}$ , and

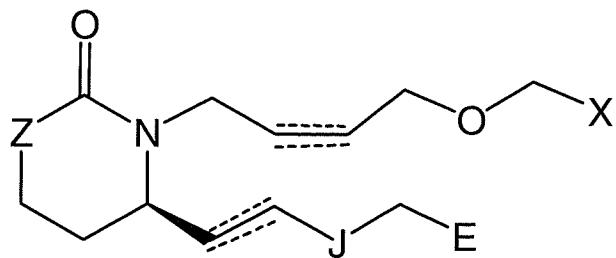


J is C=O or CHO;

R is independently H,  $\text{C}_1\text{-C}_6$  alkyl, phenyl, or biphenyl; and

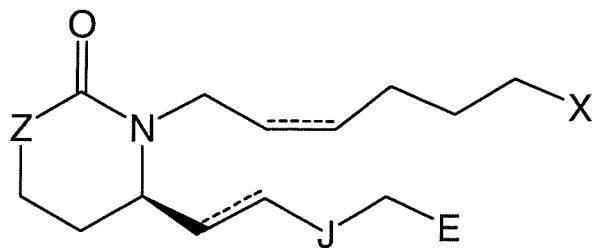
E is  $\text{C}_3\text{-C}_6$  alkyl,  $\text{C}_4\text{-C}_{10}$  cycloalkyl, phenyl or napthyl having from 0 to 2 substituents, or a heteroaromatic moiety having from 0 to 2 substituents, wherein said substituents comprise up to 4 non-hydrogen atoms.

CLAIM 2. (Currently Amended) The compound of claim 1 comprising  
represented by the formula:



or a pharmaceutically acceptable salt or a prodrug thereof.

CLAIM 3. (Currently Amended) The compound of claim 1 comprising  
represented by the formula:



or a pharmaceutically acceptable salt or a prodrug thereof.

CLAIM 4. (Cancelled)

CLAIM 5. (Cancelled)

CLAIM 6. (Cancelled)

CLAIM 7. (Original) The compound of claim 3 wherein J is C=O.

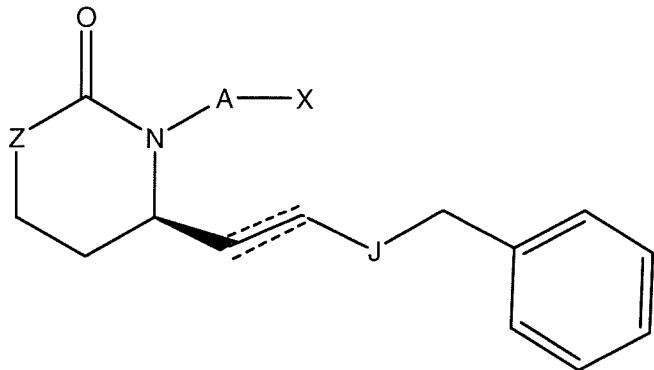
CLAIM 8. (Original) The compound of claim 3 wherein J is CHOH.

CLAIM 9. (Original) The compound of claim 3 wherein X is CO<sub>2</sub>H.

CLAIM 10. (Original) The compound of claim 3 wherein E is phenyl, thienyl, furyl, pyridinyl, napthyl, benzothienyl, or benzofuryl having from 0 to 2 substituents comprising up to 4 non-hydrogen atoms.

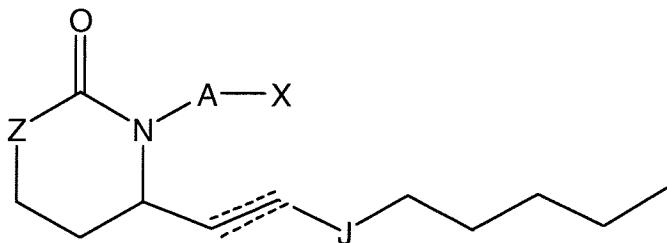
CLAIM 11. (Original) The compound of claim 3 wherein E is *n*-butyl.

CLAIM 12. (Currently Amended) The compound of claim 1 comprising represented by the formula:



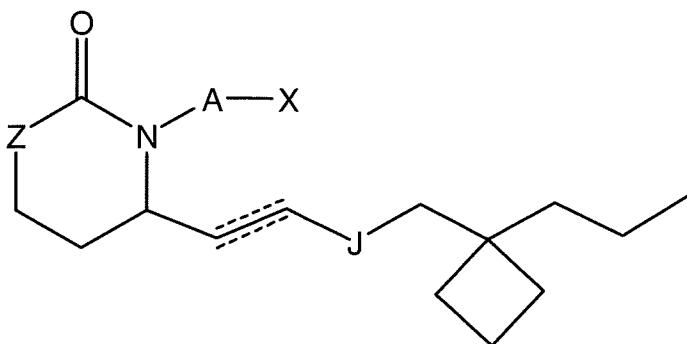
or a pharmaceutically acceptable salt or a prodrug thereof.

CLAIM 13. (Currently Amended) The compound of claim 1 comprising represented by the formula:



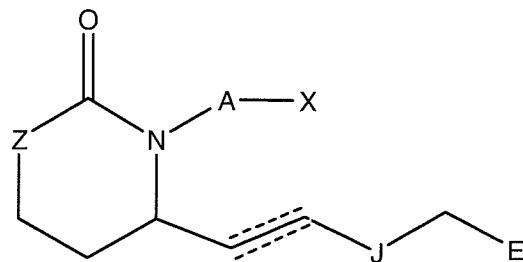
or a pharmaceutically acceptable salt or a prodrug thereof.

CLAIM 14. (Currently Amended) The compound of claim 1 comprising represented by the formula:



or a pharmaceutically acceptable salt or a prodrug thereof.

**CLAIM 15. (Currently Amended) A liquid composition comprising a compound represented by the formula:**

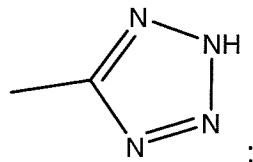


or a pharmaceutically acceptable salt or a prodrug thereof;  
wherein a dashed line represents the presence or absence of a double bond or a triple bond;

A is  $-(CH_2)_6-$ , *cis*  $-CH_2CH=CH-(CH_2)_3-$ , or  $-CH_2C\equiv C-(CH_2)_3-$ , wherein 1 or 2 carbon atoms may be substituted with S or O;

Z is O, S, or NR;

X is selected from the group consisting of  $CO_2H$ ,  $CONHR_2$ ,  $CONR_2$ ,  $CON(OR)R$ ,  $CON(CH_2CH_2OH)_2$ ,  $CONH(CH_2CH_2OH)$ ,  $CH_2OH$ ,  $P(O)(OH)_2$ ,  $CONHSO_2R$ ,  $SO_2NR_2$ ,  $SO_2NHR$ , and

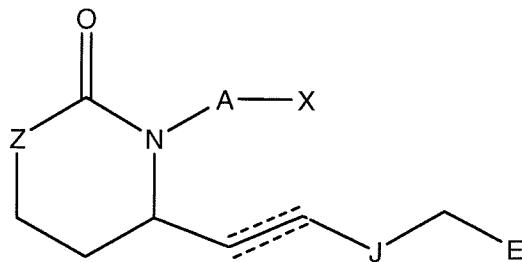


J is C=O or CHOH;

R is independently H,  $C_1-C_6$  alkyl, phenyl, or biphenyl; and

E is  $C_3-C_6$  alkyl,  $C_4-C_{10}$  cycloalkyl, phenyl or naphthyl having from 0 to 2 substituents, or a heteroaromatic moiety having from 0 to 2 substituents, wherein said substituents comprise up to 4 non-hydrogen atoms;  
wherein said liquid is formulated for ophthalmic use.

**CLAIM 16.** (Currently Amended) A method comprising administering a compound to a mammal, wherein said method is useful for the treatment of glaucoma or ocular hypertension in said mammal, said compound comprising being represented by the formula:

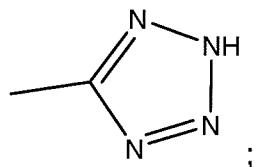


or a pharmaceutically acceptable salt or a prodrug thereof;  
wherein a dashed line represents the presence or absence of a double bond or a triple bond;

A is  $-(CH_2)_6-$ , *cis*  $-CH_2CH=CH-(CH_2)_3-$ , or  $-CH_2C\equiv C-(CH_2)_3-$ , wherein 1 or 2 carbon atoms may be substituted with S or O;

Z is O, S, or NR;

X is selected from the group consisting of  $CO_2H$ ,  $CONHR_2$ ,  $CONR_2$ ,  $CON(OR)R$ ,  $CON(CH_2CH_2OH)_2$ ,  $CONH(CH_2CH_2OH)$ ,  $CH_2OH$ ,  $P(O)(OH)_2$ ,  $CONHSO_2R$ ,  $SO_2NR_2$ ,  $SO_2NHR$ , and



J is C=O or CHOH;

R is independently H, C<sub>1</sub>-C<sub>6</sub> alkyl, phenyl, or biphenyl; and

E is C<sub>3</sub>-C<sub>6</sub> alkyl, C<sub>4</sub>-C<sub>10</sub> cycloalkyl, phenyl or napthyl having from 0 to 2 substituents, or a heteroaromatic moiety having from 0 to 2 substituents, wherein said substituents comprise up to 4 non-hydrogen atoms.